

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111325\
 Data File : BP026135.D
 Acq On : 13 Nov 2025 19:45
 Operator : RC/JU
 Sample : Q3601-09DL 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-3DL

Quant Time: Nov 14 03:54:07 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 30 11:51:34 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.672	152	326524	20.000	ng	-0.02
21) Naphthalene-d8	10.443	136	1257617	20.000	ng	-0.01
39) Acenaphthene-d10	14.313	164	758862	20.000	ng	0.00
64) Phenanthrene-d10	17.125	188	1552676	20.000	ng	0.00
76) Chrysene-d12	21.583	240	1568278	20.000	ng	0.02
86) Perylene-d12	24.907	264	1769470	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.296	112	723468	36.561	ng	0.00
7) Phenol-d6	6.855	99	914968	36.328	ng	0.00
23) Nitrobenzene-d5	8.808	82	524485	25.999	ng	-0.02
42) 2,4,6-Tribromophenol	15.831	330	344274	41.964	ng	0.00
45) 2-Fluorobiphenyl	12.919	172	1278358	24.206	ng	-0.01
79) Terphenyl-d14	19.866	244	1717705	22.253	ng	0.00
Target Compounds						
						Qvalue
52) Acenaphthene	14.378	154	95252	2.082	ng	97
58) Fluorene	15.384	166	121387	2.242	ng	98
71) Phenanthrene	17.166	178	2466541	27.293	ng	100
72) Anthracene	17.254	178	572047	6.386	ng	99
73) Carbazole	17.537	167	199582	2.354	ng	100
75) Fluoranthene	19.266	202	3727656	35.359	ng	98
78) Pyrene	19.642	202	3519801	33.184	ng	100
80) Butylbenzylphthalate	20.607	149	29172	4.082	ng	# 57
81) Benzo(a)anthracene	21.566	228	1787396	16.586	ng	96
83) Chrysene	21.630	228	1644798	16.112	ng	98
84) Bis(2-ethylhexyl)phtha...	21.530	149	55906	3.434	ng	# 95
87) Indeno(1,2,3-cd)pyrene	28.665	276	1143083	8.750	ng	# 89
88) Benzo(b)fluoranthene	23.854	252	2183000	19.873	ng	96
89) Benzo(k)fluoranthene	23.919	252	683453	6.093	ng	96
90) Benzo(a)pyrene	24.736	252	1639274	16.244	ng	96
91) Dibenzo(a,h)anthracene	28.748	278	277682	2.612	ng	# 90
92) Benzo(g,h,i)perylene	29.830	276	1091611	10.671	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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