

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031125\
 Data File : BF141918.D
 Acq On : 11 Mar 2025 16:12
 Operator : RC/JU
 Sample : Q1514-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ENV-105-SB01

Manual Integrations
 APPROVED

Reviewed By :Anahy Claudio 03/12/2025
 Supervised By :Jagrut Upadhyay 03/12/2025

Quant Time: Mar 11 16:52:37 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 10 15:46:22 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	150957	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	586367	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	300301	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	431301	20.000	ng	0.00
76) Chrysene-d12	14.039	240	338343	20.000	ng	0.00
86) Perylene-d12	15.521	264	250489	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	749092	82.819	ng	0.00
7) Phenol-d6	6.498	99	900241	78.171	ng	0.00
23) Nitrobenzene-d5	7.434	82	582366	55.892	ng	-0.01
42) 2,4,6-Tribromophenol	10.698	330	254726	66.862	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	1157452	58.617	ng	0.00
79) Terphenyl-d14	12.986	244	1029420	44.982	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	11.422	178	219318	9.414	ng	99
72) Anthracene	11.474	178	47239m	2.021	ng	
75) Fluoranthene	12.610	202	205658	8.322	ng	99
78) Pyrene	12.839	202	190783	6.519	ng	96
81) Benzo(a)anthracene	14.027	228	92270	4.156	ng	# 88
83) Chrysene	14.063	228	82029	4.076	ng	# 86
88) Benzo(b)fluoranthene	15.086	252	68107m	3.967	ng	
90) Benzo(a)pyrene	15.457	252	46734	3.479	ng	# 64

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

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