

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082919\
 Data File : BM022427.D
 Acq On : 30 Aug 2019 02:51
 Operator : HP/JU
 Sample : K4514-14
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 S301A-J-2.0-2.5-082319

Quant Time: Aug 30 08:54:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 29 16:51:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	18038	20.00	ng	0.00
21) Naphthalene-d8	10.30	136	62585	20.00	ng	0.00
39) Acenaphthene-d10	14.19	164	39491	20.00	ng	0.00
64) Phenanthrene-d10	16.95	188	92482	20.00	ng	0.00
76) Chrysene-d12	21.17	240	127513	20.00	ng	0.00
87) Perylene-d12	23.36	264	154021	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.15	112	141316	137.52	ng	0.00
7) Phenol-d6	6.72	99	181682	132.21	ng	0.00
23) Nitrobenzene-d5	8.69	82	120328	84.21	ng	0.00
42) 2,4,6-Tribromophenol	15.70	330	94573	117.03	ng	0.00
45) 2-Fluorobiphenyl	12.79	172	217388	62.36	ng	0.00
79) Terphenyl-d14	19.60	244	484075	50.10	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.75	94	5370	3.890	ng	# 74
50) Dimethylphthalate	13.66	163	35124	10.226	ng	98
71) Phenanthrene	17.00	178	14665	2.720	ng	95
75) Fluoranthene	19.03	202	41439	5.772	ng	96
78) Pyrene	19.39	202	39542	4.550	ng	# 96
81) Benzo(a)anthracene	21.15	228	25757	2.814	ng	96
83) Chrysene	21.20	228	25986	2.983	ng	99
84) Bis(2-ethylhexyl)phthalate	21.07	149	9060	2.200	ng	99
86) Indeno(1,2,3-cd)pyrene	25.54	276	17535	2.001	ng	98
88) Benzo(b)fluoranthene	22.71	252	33886	3.167	ng	# 95
90) Benzo(a)pyrene	23.26	252	24953	2.627	ng	# 98
92) Benzo(a,h,i)perylene	26.21	276	16543	2.004	ng	95

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

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