

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072922\
 Data File : BM035996.D
 Acq On : 29 Jul 2022 17:17
 Operator : CG/JU
 Sample : PB146501BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK501

Quant Time: Jul 30 00:24:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 30 00:20:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	242114	20.000	ng/u1	0.00
20) Naphthalene-d8	10.681	136	1007236	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.516	164	589121	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	1194443	20.000	ng/u1	0.00
79) Chrysene-d12	21.427	240	1104719	20.000	ng/u1	0.00
88) Perylene-d12	23.797	264	1116169	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	40511	6.296	ng/uL	0.00
4) Pyridine-d5	3.693	84	526379	29.862	ng/u1	0.00
7) Phenol-d5	7.040	99	623602	30.714	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	402021	29.702	ng/u1	0.00
11) 2-Chlorophenol-d4	7.404	132	547314	32.476	ng/u1	0.00
15) 4-Methylphenol-d8	8.587	113	485894	29.541	ng/u1	0.00
21) Nitrobenzene-d5	9.045	128	258350	32.369	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	259995	33.435	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	447119	31.792	ng/u1	0.00
31) 4-Chloroaniline-d4	10.822	131	707379	29.711	ng/u1	0.00
46) Dimethylphthalate-d6	13.927	166	1336556	33.306	ng/u1	0.00
49) Acenaphthylene-d8	14.210	160	1706511	33.329	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	221510	26.880	ng/u1	0.00
60) Fluorene-d10	15.504	176	1194169	32.931	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	153377	23.889	ng/u1	0.00
73) Anthracene-d10	17.351	188	1815938	33.628	ng/u1	0.00
81) Pyrene-d10	19.645	212	2093464	34.193	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.644	264	1973322	34.823	ng/u1	0.00

Target Compounds	Qvalue

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

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