

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102023\
 Data File : BP017738.D
 Acq On : 20 Oct 2023 11:37
 Operator : MA/JU
 Sample : PB156394BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK394

Quant Time: Oct 20 22:39:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	126888	20.000	ng/u1	0.01
20) Naphthalene-d8	10.728	136	485444	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.593	164	289409	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.387	188	599788	20.000	ng/u1	# 0.01
79) Chrysene-d12	21.533	240	491905	20.000	ng/u1	0.02
88) Perylene-d12	24.098	264	499568	20.000	ng/u1	0.03

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	25287	7.414	ng/uL	0.00
4) Pyridine-d5	3.693	84	325039	36.398	ng/u1	0.00
7) Phenol-d5	7.058	99	362333	34.315	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.246	67	243251	37.303	ng/u1	0.01
11) 2-Chlorophenol-d4	7.416	132	308857	35.863	ng/u1	0.00
15) 4-Methylphenol-d8	8.622	113	283366	34.174	ng/u1	0.01
21) Nitrobenzene-d5	9.105	128	136696	34.751	ng/u1	0.00
24) 2-Nitrophenol-d4	9.822	143	151568	34.873	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.352	165	267467	31.372	ng/u1	0.01
31) 4-Chloroaniline-d4	10.905	131	362051	32.449	ng/u1	0.01
46) Dimethylphthalate-d6	14.016	166	841112	34.621	ng/u1	0.01
49) Acenaphthylene-d8	14.293	160	925867	34.172	ng/u1	0.01
54) 4-Nitrophenol-d4	14.845	143	94883	27.409	ng/u1	0.02
60) Fluorene-d10	15.604	176	709475	33.783	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.745	200	104775	26.520	ng/u1	0.01
73) Anthracene-d10	17.492	188	1052469	34.338	ng/u1	0.02
81) Pyrene-d10	19.751	212	1225218	39.728	ng/u1	0.02
92) Benzo(a)pyrene-d12	23.927	264	1034056	37.164	ng/u1	0.03

Target Compounds	Qvalue

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

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