

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
 Data File : BP018658.D
 Acq On : 07 Dec 2023 03:40
 Operator : MA/JU
 Sample : PB157356BL
 Misc :
 ALS Vial : 96 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK356

Quant Time: Dec 07 04:55:11 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 04 21:39:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	409754	20.000	ng/u1	0.00
20) Naphthalene-d8	10.646	136	1607291	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.481	164	923930	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.228	188	1747264	20.000	ng/u1	0.00
79) Chrysene-d12	21.316	240	1634370	20.000	ng/u1	0.00
88) Perylene-d12	23.686	264	2027535	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	66054	5.493	ng/uL	0.00
4) Pyridine-d5	3.681	84	807043	25.057	ng/u1	0.00
7) Phenol-d5	7.016	99	1072996	26.094	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	646767	26.538	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	832915	25.959	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	666771	20.036	ng/u1	0.00
21) Nitrobenzene-d5	9.011	128	344684	24.095	ng/u1	0.00
24) 2-Nitrophenol-d4	9.734	143	436977	29.463	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	791045	26.084	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.787	131	849306	21.429	ng/u1	0.00
46) Dimethylphthalate-d6	13.893	166	2248488	27.267	ng/u1	-0.01
49) Acenaphthylene-d8	14.169	160	2567554	28.277	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.681	143	325430	24.042	ng/u1	0.00
60) Fluorene-d10	15.469	176	1850026	26.727	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.587	200	245046	23.390	ng/u1	0.00
73) Anthracene-d10	17.322	188	2615095	28.469	ng/u1	0.00
81) Pyrene-d10	19.563	212	2964311	23.300	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	3338104	28.236	ng/u1	0.00

Target Compounds Qvalue

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

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