

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101623\
 Data File : BP017683.D
 Acq On : 16 Oct 2023 15:39
 Operator : MA/JU
 Sample : PB156265BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK265

Quant Time: Oct 17 00:33:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 16:14:36 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	87823	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	344289	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.604	164	211322	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	489028	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	521605	20.000	ng/u1	0.00
88) Perylene-d12	24.115	264	619477	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	17646	7.408	ng/uL	0.00
4) Pyridine-d5	3.699	84	225307	34.216	ng/u1	0.00
7) Phenol-d5	7.058	99	273381	33.436	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	179502	36.201	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	230996	36.499	ng/u1	0.00
15) 4-Methylphenol-d8	8.622	113	209614	32.097	ng/u1	0.00
21) Nitrobenzene-d5	9.110	128	104447	36.118	ng/u1	0.00
24) 2-Nitrophenol-d4	9.828	143	112459	34.581	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	199116	32.577	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	304054	34.706	ng/u1	0.00
46) Dimethylphthalate-d6	14.022	166	680751	37.195	ng/u1	0.00
49) Acenaphthylene-d8	14.298	160	728344	36.956	ng/u1	0.00
54) 4-Nitrophenol-d4	14.857	143	79950	26.120	ng/u1	0.00
60) Fluorene-d10	15.610	176	598182	38.423	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	84512	24.450	ng/u1	0.00
73) Anthracene-d10	17.498	188	977300	38.712	ng/u1	0.00
81) Pyrene-d10	19.757	212	1205243	38.391	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.945	264	1345996	38.704	ng/u1	0.00

Target Compounds Qvalue

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

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