

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101623\
 Data File : BP017674.D
 Acq On : 16 Oct 2023 09:59
 Operator : MA/JU
 Sample : PB156048BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK048

Quant Time: Oct 17 00:26:13 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	7.905	152	129124	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	517453	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.604	164	324927	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	768897	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	799379	20.000	ng/u1	0.00
88) Perylene-d12	24.115	264	885759	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	20139	5.750	ng/uL	0.00
4) Pyridine-d5	3.699	84	257635	26.611	ng/u1	0.00
7) Phenol-d5	7.058	99	320885	26.693	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	211917	29.068	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	278299	29.908	ng/u1	0.00
15) 4-Methylphenol-d8	8.628	113	261029	27.185	ng/u1	0.00
21) Nitrobenzene-d5	9.116	128	128949	29.669	ng/u1	0.00
24) 2-Nitrophenol-d4	9.828	143	136713	27.971	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	235770	25.665	ng/u1	0.00
31) 4-Chloroaniline-d4	10.916	131	364539	27.685	ng/u1	0.00
46) Dimethylphthalate-d6	14.022	166	886931	31.517	ng/u1	0.00
49) Acenaphthylene-d8	14.304	160	922123	30.429	ng/u1	0.00
54) 4-Nitrophenol-d4	14.857	143	111137	23.614	ng/u1	0.00
60) Fluorene-d10	15.610	176	755829	31.575	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	123093	22.649	ng/u1	0.00
73) Anthracene-d10	17.498	188	1270742	32.014	ng/u1	0.00
81) Pyrene-d10	19.757	212	1641080	34.110	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.945	264	1712572	34.441	ng/u1	0.00

Target Compounds	Qvalue

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

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