

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017129.D
 Acq On : 12 Sep 2023 23:01
 Operator : MA/JU
 Sample : PB155287BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK287

Quant Time: Sep 13 01:30:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:27:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	316494	20.000	ng/u1	0.00
20) Naphthalene-d8	10.781	136	1407898	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	946796	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	2193263	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	2163101	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	2394566	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	46140	6.126	ng/uL	0.00
4) Pyridine-d5	3.746	84	658642	30.571	ng/u1	0.00
7) Phenol-d5	7.105	99	795576	30.914	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	491772	30.573	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	635177	30.938	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	632784	29.669	ng/u1	0.00
21) Nitrobenzene-d5	9.152	128	305528	30.991	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	339112	32.220	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	597328	29.347	ng/u1	0.00
31) 4-Chloroaniline-d4	10.946	131	885885	28.876	ng/u1	0.00
46) Dimethylphthalate-d6	14.028	166	2026450	30.077	ng/u1	0.00
49) Acenaphthylene-d8	14.316	160	2140807	28.003	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	351975	27.029	ng/u1	0.00
60) Fluorene-d10	15.610	176	1699272	29.169	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.740	200	306707	25.030	ng/u1	0.00
73) Anthracene-d10	17.481	188	2822426	30.288	ng/u1	0.00
81) Pyrene-d10	19.722	212	3544839	31.248	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	3577037	31.650	ng/u1	-0.01

Target Compounds Qvalue

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

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