

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
 Data File : BP017290.D
 Acq On : 22 Sep 2023 17:09
 Operator : MA/JU
 Sample : PB155616BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK616

Quant Time: Sep 22 18:36:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 21 23:51:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	132161	20.000	ng/u1	0.00
20) Naphthalene-d8	10.757	136	532614	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.593	164	315477	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.363	188	723073	20.000	ng/u1	0.00
79) Chrysene-d12	21.469	240	689939	20.000	ng/u1	0.00
88) Perylene-d12	23.992	264	774010	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	27205	8.455	ng/uL	0.00
4) Pyridine-d5	3.740	84	385935	42.905	ng/u1	0.00
7) Phenol-d5	7.093	99	468730	43.184	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	273482	41.750	ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	382550	44.437	ng/u1	0.00
15) 4-Methylphenol-d8	8.646	113	364928	43.244	ng/u1	-0.01
21) Nitrobenzene-d5	9.134	128	174964	45.440	ng/u1	0.00
24) 2-Nitrophenol-d4	9.846	143	193217	46.214	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	347438	43.985	ng/u1	0.00
31) 4-Chloroaniline-d4	10.922	131	494844	43.247	ng/u1	0.00
46) Dimethylphthalate-d6	14.010	166	1040206	46.027	ng/u1	0.00
49) Acenaphthylene-d8	14.293	160	1158017	44.552	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	166671	42.331	ng/u1	0.00
60) Fluorene-d10	15.592	176	905852	46.558	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	114856	28.183	ng/u1	0.00
73) Anthracene-d10	17.469	188	1542601	48.103	ng/u1	0.00
81) Pyrene-d10	19.710	212	1878006	50.335	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.827	264	1853955	49.749	ng/u1	0.00

Target Compounds	Qvalue

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

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