

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091023\
 Data File : BP017101.D
 Acq On : 11 Sep 2023 13:05
 Operator : MA/JU
 Sample : 04267-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH4F1

Quant Time: Sep 12 04:05:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 06:13:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	266992	20.000	ng/u1	0.00
20) Naphthalene-d8	10.786	136	1177848	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	714997	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	1517791	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	1151558	20.000	ng/u1	0.00
88) Perylene-d12	24.015	264	1099278	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.316	96	51751	7.398	ng/uL	0.00
4) Pyridine-d5	3.757	84	166871	7.911	ng/u1	0.00
7) Phenol-d5	7.110	99	223138	9.194	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.298	67	623734	38.319	ng/u1	0.00
11) 2-Chlorophenol-d4	7.481	132	532830	30.376	ng/u1	0.00
15) 4-Methylphenol-d8	8.669	113	385581	20.084	ng/u1	0.00
21) Nitrobenzene-d5	9.157	128	331431	39.524	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	331539	38.307	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.392	165	562799	34.294	ng/u1	0.00
31) 4-Chloroaniline-d4	10.945	131	789397	29.997	ng/u1	0.00
46) Dimethylphthalate-d6	14.033	166	2109367	41.461	ng/u1	0.00
49) Acenaphthylene-d8	14.316	160	2299087	39.687	ng/u1	0.00
54) 4-Nitrophenol-d4	14.833	143	64782	6.172	ng/u1	0.00
60) Fluorene-d10	15.616	176	1790692	41.477	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	248695	29.035	ng/u1	0.00
73) Anthracene-d10	17.486	188	2835536	43.585	ng/u1	0.00
81) Pyrene-d10	19.727	212	3091482	50.978	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.851	264	2332767	44.204	ng/u1	0.00

Target Compounds	Qvalue

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

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