

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092823\
 Data File : BP017417.D
 Acq On : 28 Sep 2023 16:57
 Operator : MA/JU
 Sample : 04417-26
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYF2

Quant Time: Sep 28 23:15:42 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 28 15:31:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	128046	20.000	ng/u1	0.02
20) Naphthalene-d8	10.816	136	509805	20.000	ng/u1	0.02
38) Acenaphthene-d10	14.669	164	302122	20.000	ng/u1	0.02
64) Phenanthrene-d10	17.457	188	633082	20.000	ng/u1	0.02
79) Chrysene-d12	21.592	240	592188	20.000	ng/u1	0.02
88) Perylene-d12	24.186	264	721936	20.000	ng/u1	0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	5307	1.702	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.128	99	86178	8.195	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.311	67	60394	9.516	ng/u1	0.01
11) 2-Chlorophenol-d4	7.493	132	71254	8.543	ng/u1	0.02
15) 4-Methylphenol-d8	8.693	113	23872	2.920	ng/u1	0.02
21) Nitrobenzene-d5	9.181	128	38486	10.442	ng/u1	0.02
24) 2-Nitrophenol-d4	9.899	143	25654	6.411	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.434	165	46829	6.194	ng/u1	0.02
31) 4-Chloroaniline-d4	10.981	131	62770	5.731	ng/u1	0.02
46) Dimethylphthalate-d6	14.081	166	230543	10.652	ng/u1	0.02
49) Acenaphthylene-d8	14.369	160	265653	10.672	ng/u1	0.02
54) 4-Nitrophenol-d4	14.887	143	144	0.038	ng/u1	0.00
60) Fluorene-d10	15.675	176	205239	11.015	ng/u1	0.02
65) 4,6-Dinitro-2-methylph...	15.798	200	101	0.028	ng/u1	0.00
73) Anthracene-d10	17.557	188	295010	10.507	ng/u1	0.02
81) Pyrene-d10	19.810	212	360252	11.249	ng/u1	0.02
92) Benzo(a)pyrene-d12	24.015	264	376211	10.823	ng/u1	0.04
Target Compounds						
					Qvalue	
16) Acetophenone	8.828	105	53231	4.213	ng/u1	97

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

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