

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010828.D
 Acq On : 25 Jun 2022 15:02
 Operator : CG/JU
 Sample : N3355-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EY0C9

Quant Time: Jun 26 22:52:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	404473	20.000	ng/u1	0.00
20) Naphthalene-d8	10.516	136	1635084	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.375	164	868967	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.140	188	1565347	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.251	240	1470596	20.000	ng/u1	# 0.00
88) Perylene-d12	23.622	264	1597307	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	31118	2.910	ng/uL	0.00
4) Pyridine-d5	3.617	84	215895	7.585	ng/u1	0.01
7) Phenol-d5	6.899	99	294600	9.201	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.058	67	736946	35.194	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	781586	29.422	ng/u1	0.00
15) 4-Methylphenol-d8	8.440	113	482406	19.059	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	448068	43.047	ng/u1	0.00
24) 2-Nitrophenol-d4	9.599	143	413377	48.003	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.140	165	690020	32.429	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	741386	21.345	ng/u1	0.00
46) Dimethylphthalate-d6	13.799	166	2233719	37.526	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	2700757	35.550	ng/u1	0.00
54) 4-Nitrophenol-d4	14.593	143	72716	7.821	ng/u1	0.01
60) Fluorene-d10	15.375	176	1892871	37.056	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.493	200	163367	31.501	ng/u1	0.00
73) Anthracene-d10	17.240	188	2724257	39.117	ng/u1	0.00
81) Pyrene-d10	19.486	212	3045103	38.007	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.469	264	3235943	40.997	ng/u1	0.00

Target Compounds	Qvalue

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

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